

* * * * * Welcome to STN International * * * * *

NEWS 1 Web Page URLs for STN Seminar Schedule - N. America
 NEWS 2 "Ask CAS" for self-help around the clock
 NEWS 3 May 12 EXTEND option available in structure searching
 NEWS 4 May 12 Polymer links for the POLYLINK command completed in REGISTRY
 NEWS 5 May 27 New UPM (Update Code Maximum) field for more efficient patent SDIs in CPlus
 NEWS 6 May 27 CPlus super roles and document types searchable in REGISTRY
 NEWS 7 Jun 22 STN Patent Forums to be held July 19-22, 2004
 NEWS 8 Jun 28 Additional enzyme-catalyzed reactions added to CASREACT
 NEWS 9 Jun 28 ANTE, AQUALINE, BIOENG, CIVILENG, ENVIROENG, MECHENG, and WATER from CSA now available on STN(R)

NEWS EXPRESS MARCH 31 CURRENT WINDOWS VERSION IS V7.00A, CURRENT MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP), AND CURRENT DISCOVER FILE IS DATED 26 APRIL 2004

NEWS HOURS STN Operating Hours Plus Help Desk Availability
 NEWS INTER General Internet Information
 NEWS LOGIN Welcome Banner and News Items
 NEWS PHONE Direct Dial and Telecommunication Network Access to STN
 NEWS WWW CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that specific topic.

All use of STN is subject to the provisions of the STN Customer agreement. Please note that this agreement limits use to scientific research. Use for software development or design or implementation of commercial gateways or other similar uses is prohibited and may result in loss of user privileges and other penalties.

* * * * * STN Columbus * * * * *

FILE 'HOME' ENTERED AT 03:56:14 ON 08 JUL 2004

=> file reg

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
0.21	0.21

FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 03:56:21 ON 08 JUL 2004

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

COPYRIGHT (C) 2004 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 6 JUL 2004 HIGHEST RN 705249-96-3

DICTIONARY FILE UPDATES: 6 JUL 2004 HIGHEST RN 705249-96-3

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 6, 2004

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See [HELP CROSSOVER](#) for details.

Experimental and calculated property data are now available. For more

information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at:
<http://www.cas.org/ONLINE/DBSS/registryss.html>

=>

L1 STRUCTURE UPLOADED

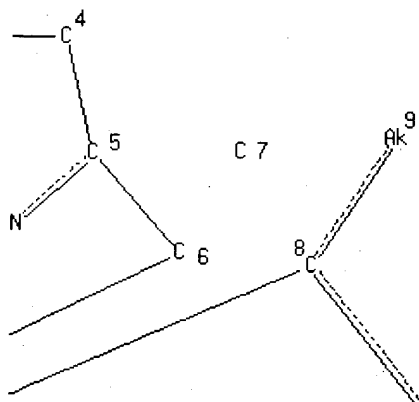
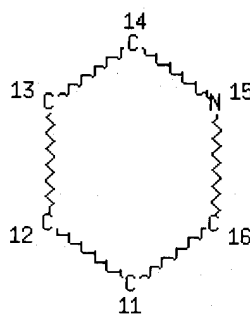
=> d l1

L1 HAS NO ANSWERS

L1 STR



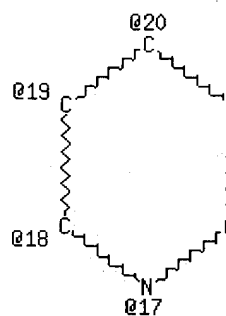
Page 1-A



Page 1-B

G20
23
Page 2-A

AK
Q10



Page 2-B
Q21

Q22
Page 2-C
REP G20=(0-2) 7-6 7-8
VPA 10-17/18/19/20/21/22 S
NODE ATTRIBUTES:

NSPEC	IS R	AT	1
NSPEC	IS R	AT	2
NSPEC	IS R	AT	3
NSPEC	IS R	AT	4
NSPEC	IS R	AT	5
NSPEC	IS C	AT	6
NSPEC	IS C	AT	7
NSPEC	IS C	AT	8
NSPEC	IS C	AT	9
NSPEC	IS C	AT	10
NSPEC	IS R	AT	11
NSPEC	IS R	AT	12
NSPEC	IS R	AT	13
NSPEC	IS R	AT	14

NSPEC IS R AT 15
 NSPEC IS R AT 16
 NSPEC IS R AT 17
 NSPEC IS R AT 18
 NSPEC IS R AT 19
 NSPEC IS R AT 20
 NSPEC IS R AT 21
 NSPEC IS R AT 22
 NSPEC IS C AT 23
 DEFAULT MLEVEL IS ATOM
 MLEVEL IS CLASS AT 6 7 8 9 10
 DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:

RSPEC I

NUMBER OF NODES IS 23

STEREO ATTRIBUTES: NONE

=> s ll

SAMPLE SEARCH INITIATED 04:01:06 FILE 'REGISTRY'
 SAMPLE SCREEN SEARCH COMPLETED - 6135 TO ITERATE

16.3% PROCESSED 1000 ITERATIONS
 INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)
 SEARCH TIME: 00.00.01

0 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
 BATCH **COMPLETE**
 PROJECTED ITERATIONS: 118005 TO 127395
 PROJECTED ANSWERS: 0 TO 0

L2 0 SEA SSS SAM L1

=> s ll full

THE ESTIMATED SEARCH COST FOR FILE 'REGISTRY' IS 155.00 U.S. DOLLARS
 DO YOU WANT TO CONTINUE WITH THIS REQUEST? (Y)/N or END:y
 FULL SEARCH INITIATED 04:01:11 FILE 'REGISTRY'
 FULL SCREEN SEARCH COMPLETED - 124961 TO ITERATE

100.0% PROCESSED 124961 ITERATIONS
 SEARCH TIME: 00.00.08

4 ANSWERS

L3 4 SEA SSS FUL L1

=> file hcaplus

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
158.36	158.57

FULL ESTIMATED COST

FILE 'HCAPLUS' ENTERED AT 04:01:22 ON 08 JUL 2004
 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.
 PLEASE SEE "HELP USAGETERMS" FOR DETAILS.
 COPYRIGHT (C) 2004 AMERICAN CHEMICAL SOCIETY (ACS)

Copyright of the articles to which records in this database refer is held by the publishers listed in the PUBLISHER (PB) field (available for records published or updated in Chemical Abstracts after December

26, 1996), unless otherwise indicated in the original publications. The CA Lexicon is the copyrighted intellectual property of the the American Chemical Society and is provided to assist you in searching databases on STN. Any dissemination, distribution, copying, or storing of this information, without the prior written consent of CAS, is strictly prohibited.

FILE COVERS 1907 - 8 Jul 2004 VOL 141 ISS 2
FILE LAST UPDATED: 7 Jul 2004 (20040707/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> s l3

L4 2 L3

=> s l4 and mullican, m?/au

57 MULLICAN, M?/AU

L5 0 L4 AND MULLICAN, M?/AU

=> s l4 and lauffer, d?/au

46 LAUFFER, D?/AU

L6 0 L4 AND LAUFFER, D?/AU

=> s l4 and tung, r?/au

302 TUNG, R?/AU

L7 0 L4 AND TUNG, R?/AU

=> d l4, ibib abs fhitr, 1-2

L4 ANSWER 1 OF 2 HCAPLUS COPYRIGHT 2004 ACS on STN

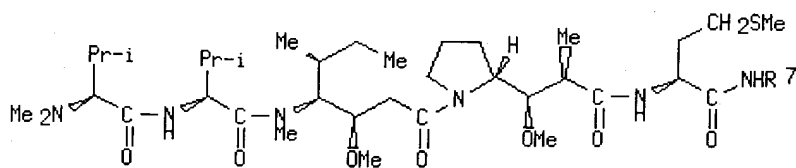
Full Text	Citing References
-----------	-------------------

ACCESSION NUMBER:	2001:185778 HCAPLUS
DOCUMENT NUMBER:	134:237837
TITLE:	Preparation of dolastatin peptides
INVENTOR(S):	Petit, George R.; Srirangam, Jayaram K.; Williams, Michael D.; Durkin, Kieran P. M.; Barlozzari, Teresa; Kling, Andreas; Janssen, Bernd; Haupt, Andreas
PATENT ASSIGNEE(S):	Basf Aktiengesellschaft, Germany; Arizona Board of Regents
SOURCE:	PCT Int. Appl., 55 pp. CODEN: PIXXD2
DOCUMENT TYPE:	Patent
LANGUAGE:	English
FAMILY ACC. NUM. COUNT:	1
<u>PATENT INFORMATION:</u>	

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2001018032	A2	20010315	WO 2000-US24658	20000908
WO 2001018032	A3	20020711		
W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CR, CU, CZ, DE, DK, DM, DZ, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY,				

DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, BF, BJ,
CF, CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG

US 6323315 B1 20011127 US 2000-539935 20000331
PRIORITY APPLN. INFO.: US 1999-394962 A1 19990910
US 2000-539935 A1 20000331
OTHER SOURCE(S): MARPAT 134:237837
GI



AB Peptides R1R2NCHR3CONHCHR4CONMeCHR5CH(OMe)CH2CONRCHRCH(OMe)CHMeCO-An-NR6R7
[R2 = (CH2)3; R1-R5 are each independently H or C1-C6 alkyl; A is a
methionyl, phenylalanyl or phenylglycyl residue; n is 0 or 1; R6 is H and
R7 is a carbocyclic, arom., alkyl, pyridylalkyl, or heterocyclic group or
R6 is benzyl or carbalkoxy and R7 is 2-thiazolyl] or their
pharmaceutically acceptable salts were prepd. for use as cell growth
inhibitors. Thus, I [R7 = bicyclo[3.3.0]octan-1-yl] was prepd. by soln.
phase methods and evaluated for in vitro cytotoxicity against a panel of
cultured cancer cell lines, including OVCAR-3 (ovarian cancer), ED50 = 3.1
x 10⁻⁴ µg/mL.

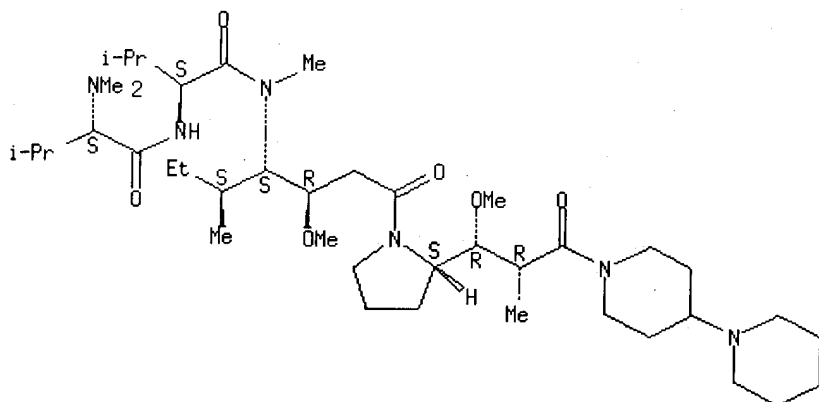
IT 329792-17-8P

RL: BAC (Biological activity or effector, except adverse); BSU (Biological
study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use);
BIOL (Biological study); PREP (Preparation); USES (Uses)
(prepn. of dolastatin peptides)

RN 329792-17-8 HCAPLUS

CN L-Valinamide, N,N-dimethyl-L-valyl-N-[(1S,2R)-4-[(2S)-2-[(1R,2R)-3-[1,4'-
bipiperidin]-1'-yl]-1-methoxy-2-methyl-3-oxopropyl]-1-pyrrolidinyl]-2-
methoxy-1-[(1S)-1-methylpropyl]-4-oxobutyl]-N-methyl- (9CI) (CA INDEX
NAME)

Absolute stereochemistry. Rotation (-).

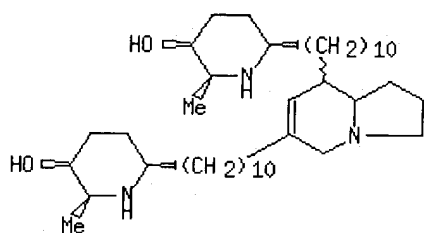


L4 ANSWER 2 OF 2 HCAPLUS COPYRIGHT 2004 ACS on STN

Full Text Citing
References

ACCESSION NUMBER: 1981:44042 HCAPLUS
DOCUMENT NUMBER: 94:44042

TITLE: Organic natural products. Part 176. Structure of the alkaloid juliprosopine from *Prosopis juliflora* A. DC
 AUTHOR(S): Ott-Longoni, Rita; Viswanathan, Narayanaier; Hesse, Manfred
 CORPORATE SOURCE: Org.-Chem. Inst., Univ. Zurich, Zurich, CH-8057, Switz.
 SOURCE: Helvetica Chimica Acta (1980), 63(7), 2119-29
 CODEN: HCACAV; ISSN: 0018-019X
 DOCUMENT TYPE: Journal
 LANGUAGE: German
 GI



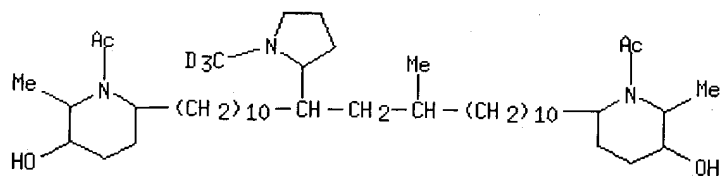
AB A new alkaloid, juliprosopine (I) was isolated from *P. juliflora* and its structure elucidated. The piperidine moiety could be elucidated by spectroscopic data (mass spectroscopy, ^1H -NMR, ^{13}C -NMR) of I and various derivs. The central hexahydroindolizine part and its substitution pattern were detd. mainly by the Hofmann-degrdn. product. I is built up in nature by 2 monomeric *Prosopis* alkaloids and 1 dihydropyrrol unit.

IT 76202-04-5P

RL: SPN (Synthetic preparation); PREP (Preparation)
 (prepn. of)

RN 76202-04-5 HCAPLUS

CN 3-Piperidinol, 6,6'-[11-methyl-13-[1-(methyl- d_3)-2-pyrrolidinyl]-1,23-tricosanediyl]bis[1-acetyl-2-methyl- (9CI) (CA INDEX NAME)



=> file caold

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
17.42	175.99

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION
-1.47	-1.47

CA SUBSCRIBER PRICE

FILE 'CAOLD' ENTERED AT 04:03:14 ON 08 JUL 2004

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

COPYRIGHT (C) 2004 AMERICAN CHEMICAL SOCIETY (ACS)

FILE COVERS 1907-1966
 FILE LAST UPDATED: 01 May 1997 (19970501/UP)

This file contains CAS Registry Numbers for easy and accurate substance identification. Title keywords, authors, patent assignees, and patent information, e.g., patent numbers, are now searchable from 1907-1966. TIFF images of CA abstracts printed between 1907-1966 are available in the PAGE display formats.

This file supports REGISTRY for direct browsing and searching of all substance data from the REGISTRY file. Enter HELP FIRST for more information.

=> d his

(FILE 'HOME' ENTERED AT 03:56:14 ON 08 JUL 2004)

FILE 'REGISTRY' ENTERED AT 03:56:21 ON 08 JUL 2004

L1 STRUCTURE UPLOADED
 L2 0 S L1
 L3 4 S L1 FULL

FILE 'HCAPLUS' ENTERED AT 04:01:22 ON 08 JUL 2004

L4 2 S L3
 L5 0 S L4 AND MULLICAN, M?/AU
 L6 0 S L4 AND LAUFFER, D?/AU
 L7 0 S L4 AND TUNG, R?/AU

FILE 'CAOLD' ENTERED AT 04:03:14 ON 08 JUL 2004

=> s l3

L8 0 L3

=> file bielstein

'BIELSTEIN' IS NOT A VALID FILE NAME

SESSION CONTINUES IN FILE 'CAOLD'

Enter "HELP FILE NAMES" at an arrow prompt (=>) for a list of files that are available. If you have requested multiple files, you can specify a corrected file name or you can enter "IGNORE" to continue accessing the remaining file names entered.

=> file beilstein

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
0.42	176.41

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION
0.00	-1.47

CA SUBSCRIBER PRICE

FILE 'BEILSTEIN' ENTERED AT 04:03:27 ON 08 JUL 2004

COPYRIGHT (c) 2004 Beilstein-Institut zur Foerderung der Chemischen Wissenschaften
 licensed to Beilstein GmbH and MDL Information Systems GmbH

FILE RELOADED ON OCTOBER 20, 2002

FILE LAST UPDATED ON JUNE 15, 2004

FILE COVERS 1771 TO 2003.

FILE CONTAINS 8,997,153 SUBSTANCES

>>> PLEASE NOTE: Reaction data and substance data are stored in separate documents and can not be searched together in one query.

Reaction data for BEILSTEIN compounds may be displayed immediately with the display codes PRE (preparations) and REA (reactions). A substance answer set retrieved after the search for a chemical name, a molecular formula or a structure search for example can be restricted to compounds with available reaction information by concatenation with PRE/FA, REA/FA or more general with RX/FA. The BEILSTEIN Registry Number (BRN) is the link between a BEILSTEIN compound and belonging reactions. For more detailed reaction searches BRNs can be selected from substance answer sets and searched in the next step as reaction partner BRNs - Reactant (RX.RBRN) or Product BRN (RX.PBRN). After a search for reaction details substance documents associated with reactants or products may be retrieved by searching RX.PBRNs or RX.RBRNs as BRNs. <<<

>>> FOR SEARCHING PREPARATIONS SEE HELP PRE <<<

```
*****
* PLEASE NOTE THAT THERE ARE NO FORMATS FREE OF COST.          *
* SET NOTICE FEATURE: THE COST ESTIMATES CALCULATED FOR SET NOTICE *
* ARE BASED ON THE HIGHEST PRICE CATEGORY. THEREFORE; THESE    *
* ESTIMATES MAY NOT REFLECT THE ACTUAL COSTS.                  *
* FOR PRICE INFORMATION SEE HELP COST                          *
*****
```

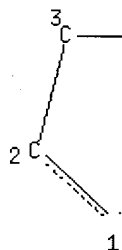
=>

L9 STRUCTURE UPLOADED

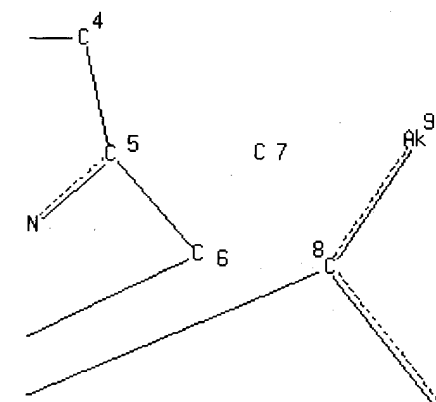
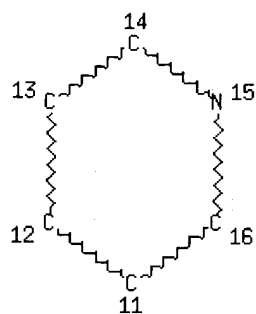
=> d 19

L9 HAS NO ANSWERS

L9 STR

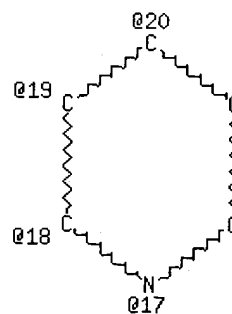


Page 1-A



Page 1-B

G20
23
Page 2-A



Page 2-B

021

022

Page 2-C

REP G20=(0-2) 7-6 7-8

VPA 10-17/18/19/20/21/22 S

NODE ATTRIBUTES:

NSPEC	IS R	AT	1
NSPEC	IS R	AT	2
NSPEC	IS R	AT	3
NSPEC	IS R	AT	4
NSPEC	IS R	AT	5
NSPEC	IS C	AT	6
NSPEC	IS C	AT	7
NSPEC	IS C	AT	8
NSPEC	IS C	AT	9
NSPEC	IS C	AT	10
NSPEC	IS R	AT	11
NSPEC	IS R	AT	12
NSPEC	IS R	AT	13
NSPEC	IS R	AT	14
NSPEC	IS R	AT	15
NSPEC	IS R	AT	16
NSPEC	IS R	AT	17
NSPEC	IS R	AT	18
NSPEC	IS R	AT	19
NSPEC	IS R	AT	20
NSPEC	IS R	AT	21
NSPEC	IS R	AT	22
NSPEC	IS C	AT	23

DEFAULT MLEVEL IS ATOM
 MLEVEL IS CLASS AT 6 7 8 9 10
 DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:

RSPEC I

NUMBER OF NODES IS 23

STEREO ATTRIBUTES: NONE

=> s 19

SAMPLE SEARCH INITIATED 04:03:45 FILE 'BEILSTEIN'
 SAMPLE SCREEN SEARCH COMPLETED - 285 TO ITERATE

100.0% PROCESSED 285 ITERATIONS
 SEARCH TIME: 00.00.04

0 ANSWERS

FULL FILE PROJECTIONS: ONLINE **COMPLETE**
 BATCH **COMPLETE**

PROJECTED ITERATIONS: 4688 TO 6712
 PROJECTED ANSWERS: 0 TO 0

L10 0 SEA SSS SAM L9

=>